

*AERONAUTICAL SYSTEMS CENTER
MAJOR SHARED RESOURCE CENTER*



*HP XC OPTERON
USER'S GUIDE*

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1. Introduction

This document provides an overview and introduction to using the HP XC Opteron pioneer system, located at the Aeronautical Systems Center (ASC) Major Shared Resource Center (MSRC). The ASC MSRC is located at Wright-Patterson Air Force Base, near Dayton, Ohio. This guide is intended to provide information so that customers who are familiar with the UNIX operating system can create and run their own programs, as well as use existing application software on the ASC MSRC HP XC Opteron system.

1.1 Assumed Background of the Reader

It is assumed that the reader of this guide has a firm grasp of the concepts required to use the UNIX operating system and to program in either the C, C++, FORTRAN 77, FORTRAN 90, or FORTRAN 95 languages. It is also assumed that the reader has read the *ASC MSRC User's Guide* (<http://www.asc.hpc.mil/customer/userdocs/ascuserg/ascuserg.pdf>), which contains site specific information about the ASC MSRC. The *ASC MSRC User's Guide* is available from the ASC MSRC Service Center at 1-888-MSRC-ASC (1-888-677-2272), (937)255-0194, or DSN 785-0194. It is also available in PostScript and PDF formats as described below in Section 1.7.

1.2 Hardware Overview

The HP XC Opteron system is a Distributed Memory system with 2 Central Processing Units (CPUs) per node. Each CPU is a 2.8 GigaHertz (GHz) AMD Opteron processor and has a 64 kilobyte (KB) Level 1 cache and 1 MB of Level 2 cache. Due to Linux Kernel and MPI memory management, users have access to a minimum of 2.25 to 2.5 GB of memory per node, depending on the number of MPI processes the user spawns.

1.3 Accessing the System

The HP XC Opteron has a total of 2048 CPUs, divided into 1024 nodes with 2 CPUs per node with 4 interactive nodes having 4 CPUs each.. The fully qualified hostname of the interactive system is *falcon.asc.hpc.mil*.

Users are only permitted to login onto the interactive nodes of the system. The other nodes are for batch jobs only. Users submit their jobs on the front-end node and the batch system will automatically start their jobs on the other systems based on the load of the system. See the *ASC MSRC User's Guide* for instructions on accessing the machines.

1.4 ASC MSRC Connectivity

Since the HP XC Opteron is an integrated component of the ASC MSRC, user files are Network File System (NFS) mounted from the ASC MSRC High Availability File Server (HAFS) system to the system. When users log into a system, their home (\$HOME) directory (which will be the current directory immediately after logging in) physically resides on the file server, but appears to be local to the system. The ASC MSRC also supplies archival storage and visualization capabilities.

1.5 ASC MSRC Startup Files

All users are provided a `.cshrc` and `.login` file in their home directory. These files reference standard setup files, maintained by the site administrators in a central

location, which set up a standard environment for all MSRC users. These files **should not** be modified.

To set up specific information for your HP XC Opteron session, such as environment variables, path information, terminal information, or command aliases, place the appropriate commands and information into files called `.personal.cshrc` and `.personal.login`. The standard startup files check your home directory for the existence of these files and executes them if found. Commands related to aliases, prompts, and some environment variables should go into `.personal.cshrc`, while commands related to the type of terminal you are using should go into `.personal.login`. See Section 3 of this guide for more details on the computing environment and the *ASC MSRC User's Guide* for more details on startup files.

1.6 The Archive Command

The archive system (\$ARC) is mounted on the interactive node of the system, however, to improve transfer speeds for batch jobs, \$ARC **is not** NFS-mounted to the batch nodes. To transfer files from \$ARC, users will have to use un-kerberized `rcp` or the archive command.

The **archive** command is a recently added tool to the ASC MSRC to help users with transferring files to and from \$ARC. The basic syntax for the archive command is:

```
archive get [getopts] file1 [file2 ...]  
archive put [putopts] file1 [file2 ...]
```

More information on the **archive** command can be found online via the *archive* man page (*man archive*).

1.7 Additional Information

Much of the information presented in this document is available online through the man pages and is accessible by typing:

```
man {command name}
```

when logged into *falcon*.

The *ASC MSRC User's Guide* and this document are all available in PostScript and PDF format. They may be downloaded via the ASC MSRC website at

<http://www.asc.hpc.mil/customer/userdocs>

2. ASC MSRC HP XC Opteron

This section details the hardware and software available on the HP XC Opteron and how they are currently configured.

2.1 Hardware

The HP XC Opteron has a total of 2048 CPUs available in batch. Each CPU is a 2.8 GHz AMD Opteron processor with a peak speed of 5.7 GFLOPS, providing a total capacity of approximately 11.64 TFLOPs. Each CPU has a primary data cache of 64 KB, a primary instruction cache of 64 KB and access to 2 GB of system memory. At this time, the system is divided into 1024 separate nodes. The system has a Distributed MultiProcessing architecture with 2048 CPUs, 4 GB of memory per node and a total of 96TB of disk space. Due to Linux Kernel and MPI memory management, users have access to a minimum of 2.25 to 2.5 GB of memory per node, depending on the number of MPI processes the user spawns.

When logging in to *falcon*, the user will be placed onto one of 4 interactive nodes due to load-balancing. Each of these nodes have identical hardware specifications to the batch nodes, however, there are 4 CPUs per node on these interactive nodes as opposed to only 2 CPUs on the batch nodes and the CPUs on the interactive nodes are 2.4 GHz as opposed to the 2.8 GHz CPUs on the batch nodes.

2.2 File System Overview

Diskspace is subdivided into several areas:

- System space (i.e., /usr, /opt.)
- /workspace

2.2.1 Workspace

Workspace is a filesystem local to each machine that batch jobs are required to run their jobs in. Workspace on the HP XC Opteron totals 96 TB.

I/O on the /workspace filesystem is quicker than I/O on a file system mounted from the network (such as \$HOME or \$ARC). /workspace is intended for the **temporary** storage of data files needed for your application. This includes (but is not limited to) grid files, restart files, input files, and output files. \$WRK is to be used rather than the /tmp and /usr/tmp directory areas to prevent possible system crashes.

There is no quota on the amount of disk space you may use in /workspace, but a file scrubber is used to automatically remove old files to prevent it from becoming filled. The current policy for removing files is on the ASC MSRC web page (http://www.asc.hpc.mil/overall/policy_procedure/policies/wrkspace_pol.php) and is subject to change based on periodic reviews.

The \$WRK file system is NOT backed up. In the event of deletion or catastrophic media failure, files and data structures are lost. It is your responsibility to transfer files that need to be saved to a location that allows permanent storage such as \$HOME or, preferably, \$ARC.

2.3 Operating System

The HP XC Opteron runs a version of Red Hat Enterprise Linux AS release 4.

2.4 Available Software

Software currently available on the HP XC Opteron includes: the PGI and GNU FORTRAN, C, and C++ compilers; MPI and many third party software packages.

3. Program Development

Program development in the HP XC Opteron computing environment is similar to that used in a typical UNIX environment. However, the user must take additional steps to utilize the multiple processors available.

3.1 Development Tools

The ASC MSRC offers many tools to help users who write their own code to develop, compile and debug their software.

The HP XC Opteron implements the PGI® High-Performance Fortran, C and C++ compilers, along with the GNU Fortran, C and C++ compilers.

The AMD Core Math Library (ACML) is installed, which provides a collection of mathematical and scientific libraries including Basic Linear Algebra Subprograms (BLAS) levels 1, 2, and 3; LAPACK; Fast Fourier Transforms (FFTs); and convolutions.

To aid users in debugging their software applications, the ASC MSRC provides the Totalview Debugger, PGI Debuggr (*pgdbg*) and the GNU Debugger (*gdb*). To launch the Totalview debugger or view it's manpage, users will have to load the Totalview module, via the *module load totalview/default* command. Once the totalview module is loaded, the command to launch Totalview is *totalview*.

The Totalview Debugger and GNU Debugger are symbolic source code debuggers that debug programs compiled by the PGI® C/C++ Compiler, the PGI® Fortran Compiler, and the GNU compilers (gcc, g++). For full source-level debugging, compile the source code with the compiler option that includes the symbol table information in the compiled executable file.

Documentation for these compilers, libraries, and tools is available online in the man page by executing a man on *pgf77*, *pgf90*, *pgf95*, *pgcc*, *pgCC*, *gcc*, *g++*, *g77*, *acml*, *totalview* (after loading the Totalview module), *pgdbg*, *gdb* or *idb*

3.2 Parallel Processing

Users may utilize multiple CPUs to execute their programs. The compilers are capable of creating parallel programs through the use of compiler directives and parallel standards such as Message Passing Interface (MPI) and OpenMP.

3.2.1 MPI

The goal of MPI is to develop a widely used standard for writing message-passing programs. As such, the interface attempts to establish a practical, portable, efficient, and flexible standard for message passing.

Users can compile MPI programs on the HP XC Opteron using the *mpicc*, *mpiCC*, *mpif77* and *mpif90* commands. These commands are wrappers around the PGI compilers that automatically link the appropriate MPI libraries for users.

For MPI FORTRAN codes:

```
mpif77 -o prog prog.f  
mpif90 -o prog prog.f
```


mpif95 -o prog prog.f

For MPI C/C++ codes:

mpicc -o prog prog.c
mpiCC -o prog prog.c

Please consult the man pages for the compiler for more information.

To execute MPI code on the HP XC Opteron users must use the ***mpirun*** command.

When using in batch, the users must use this ***mpirun*** command:

mpirun -srun -n {# of CPUs} {My Program}

When using interactively, the users must use this ***mpirun*** command:

mpirun -np {# of CPUs} {My Program}

More information on MPI can be obtained from:

<http://www.mpi-forum.org>

3.2.2 OpenMP

OpenMP is a specification for a set of compiler directives, library routines, and environment variables that can be used to specify shared memory parallelism in FORTRAN and C/C++ programs.

Creating an OpenMP program is done through OpenMP directives in the source code and by adding the -mp flag to your compile string using the PGI compilers. OpenMP is currently unsupported in the GNU compilers.

To run an OpenMP program, users must first tell the program how many threads (processors) to use. This is achieved through the OMP_NUM_THREADS environment variable. To set this variable, use the following command in *ssh*:

setenv OMP_NUM_THREADS x

where *x* is the number of CPUs you wish to utilize.

For more information on OpenMP and its directives, please see the following page:

<http://www.openmp.org>

3.2.3 MPICH

MPICH is a freely available, portable implementation of MPI, the Standard for message-passing libraries and the HP XC Opteron supports MPICH 1.2.

Users can create MPICH applications using the ***mpicc.mpich***, ***mpiCC.mpich***, ***mpif77.mpich***, and ***mpif90.mpich***.

These commands are wrappers around the PGI compilers that automatically link the appropriate MPICH libraries for users.

For MPICH FORTRAN codes:

mpif77.mpich -o prog prog.f

mpif90.mpich -o prog prog.f

For MPICH C/C++ codes:

mpicc.mpich -o prog prog.c

mpiCC.mpich -o prog prog.c

Please consult the man pages for the compiler for more information.

To execute MPICH application on the HP XC Opteron users must use the ***mpirun.mpich*** command.

When using in batch, the users must use this ***mpirun.mpich*** command:

mpirun.mpich -srun -n {# of CPUs} {My Program}

When using interactively, the users must use this ***mpirun.mpich*** command:

mpirun.mpich -np {# of CPUs} {My Program}

More information on MPI can be obtained from:

<http://www.mpi-forum.org>

3.2.4 MPI/OpenMP Hybrid Code

Some users are experimenting with codes implementing a hybrid of MPI and OpenMP calls. While the HP XC Opteron can run such code, there are some issues and limitations with doing so.

Because of the system configuration, hybrid applications are only able to run a two OpenMP processes and a single MPI process per node, up to 1024 nodes per job.

The syntax for the HP XC Opteron's mpirun command is similar to other platforms users are used to running on. Users must also set the OMP_NUM_THREADS variable, much like in the OpenMP section above.

When using in batch, the users must use this ***MPI/OpenMP*** command

setenv OMP_NUM_THREADS {Number of threads}

mpirun -srun -n {number of cpus} {My program}

When using interactively, the users must use this ***MPI/OpenMP*** command

setenv OMP_NUM_THREADS {Number of threads}

mpirun -np {number of cpus} {My program}

3.3 FORTRAN Programming

The default FORTRAN compiler on the HP XC Opteron is the PGI FORTRAN compiler. The FORTRAN compiler commands are ***pgf77***, ***pgf90*** and ***pgf95***. These optimizing and parallelizing compilers can generate 64-bit and 32-bit code. Compiling a FORTRAN program on the HP XC Opteron is similar to compiling a program on a typical UNIX system.

pgf77 -o prog prog.f

pgf90 -o prog prog.f90

prg95 -o prog prog.f95

g77 -o prog prog.f

This command creates an executable called **prog**. The program is run by typing the program name at the system prompt.

./prog

Further optimization is available through the use of compiler flags and compiler directives. Please check the **pgf77**, **pgf90**, **pgf95** and **g77** man pages for more details.

3.4 C/C++ Programming

The PGI and GNU C and C++ compilers are available on the HP XC Opteron. These compilers are capable of optimizing and parallelizing code. Compiling a C program on the HP XC Opteron is similar to compiling a C program on a typical UNIX system.

pgcc -o prog prog.c

gcc -o prog prog.c

Compiling a C++ program is also just as similar.

pgCC -o prog prog.cpp

g++ -o prog prog.cpp

These commands will create an executable program in a file called **prog**. The program is executed by entering

./prog

Further optimization is available through the use of compiler flags and compiler directives. Please consult the **icc** or **icpc** man pages for more details.

3.5 Libraries

3.5.1 Math and Science Libraries

The HP XC Opteron has ACML, a collection of mathematical and scientific libraries including BLAS levels 1, 2, and 3; LAPACK, a collection of solvers for dense linear algebra problems, including linear equations, linear least squares problems, eigenvalue problems, and singular value decomposition problems; Fast Fourier Transforms (FFTs); and convolutions. Both single-threaded and multi-threaded routines are available and select routines have been highly optimized to greatly improve performance. Users should use these library routines whenever possible.

This library is not automatically included in the link path. The user must specify the library when linking as in the following examples.

pgf77 -o prog prog.f -L/opt/acml3.0/pgi64/lib -lacml

pgcc -o prog prog.c -L/opt/acml3.0/pgi64/lib -lacml

pgCC -o prog prog.c -L/opt/acml3.0/pgi64/lib -lacml

Please consult the *ACML* web page (<http://developer.amd.com/acml.aspx>) or /opt/adml3.0/Doc on *falcon* for more details.

4. Running Jobs

4.1 Interactive Use

Interactive use is allowed, particularly for program development, including debugging and performance improvement, job preparation, job submission, and the preprocessing and postprocessing of data. Four nodes on the system are available for interactive use and interactive jobs are limited to 4 CPUs with 15 minutes of CPU time per process. Jobs with larger resource requirements must be submitted to the batch queues.

4.2 Batch Use

Load Sharing Facility (LSF), a networked subsystem for submitting, monitoring, and controlling a work load of batch jobs on one or more systems, is the batch system for the system. It provides services to monitor queue activity and to delete queued or running jobs. In the event of an orderly system shutdown, LSF jobs will be rerun from the beginning of the job (unless they are specifically marked not to be rerun). More information about LSF is available at

<http://www.platform.com/products/HPC/>

To allow users to run longer in the queues, a 2 week (336 hour) queue has been implemented, however, to keep the queue structure fair to all users, several restrictions have been put into place:

- The primary resource to schedule jobs by is the cpu hour (CPH). This quantity is equal to the wall time requested multiplied by the number of cpus and cannot exceed a value of 50,000 per user.
 $(ncpus * walltime) = CPH$

Table 1: Example CPH Values

# of Jobs	# of CPUs	Walltime	Total CPH
1	1000	48	48,000
2	256	48	24,576
25	8	96	19,200

- A user can use, at most, 1000 CPUs for MPI jobs, and 2 CPUs for SMP and MPI/OpenMP hybrid jobs.
- The maximum wall time of any queue shall not exceed 168 hours.
- A background job cannot start if there is a foreground job in any queue.

The list of queues and the upper limits of job resources for these queues are available on the web at

http://www.asc.hpc.mil/overall/policy_procedure/policies/use_policy.php

These limits are subject to change based on periodic review of system utilization and system configuration.

TIP: Because of CPH, it is not recommended that users accept the queue default walltime. If your job requires less time to run than the queue default, requesting the smaller of the two will result in a lower CPH and will allow you to **run more jobs** and **reduce your queue wait time**.

Example: User Joe submits a 4 CPU job using the queue default walltime of 336 hours. This results in a CPH of 1344. User Bob submits the same 4 CPU job, but only requests the 120 hours he needs to finish the job. His CPH would be 480. This allows user Bob to submit twice (2.8 to be exact) as many jobs as Joe in the same amount of CPH. If user Bob can alter his walltime and fit his jobs within 110 hours, he can submit 3 times the number of jobs as Joe. This also has a domino effect on the queue structure, resulting in faster throughput throughout the entire system.

4.2.1 Queueing Structure

The ASC MSRC HP XC Opteron has three queues: debug, standard, and background. These queues are available 24 hours a day, 7 days a week.

The debug queue accepts jobs that require up to 32 CPUs, 1 hour of CPU time, and 32 GB of memory. This queue is intended for short runs.

The standard queue is available for production work. Jobs that are submitted without any queue specified will go to the regular queue. Once submitted to this queue, jobs are assigned priority based upon project status and wait time in the queue.

The background queue is also available for production work. Jobs run in the background queue are not charged against a user's allocation. However, jobs in the background queue are only started when utilization of the machine is low and never when foreground jobs are waiting.

The list of queues and the upper limits of job resources for these queues are available on the web at

http://www.asc.hpc.mil/overall/policy_procedure/policies/batchqueue.php

These limits are subject to change based on periodic review of system utilization and system configuration.

4.2.2 Preparing Jobs

Before a user submits a job, they should prepare a job script. A job script is a UNIX shell script that contains all the commands the user will execute during the job. LSF will place the error and output files in the directory the job was submitted from, so scripts must be written with this in mind. Here is a sample job script.

```

#
#Change to WORK_DIR directory and copy input file.
#
cd $WORK_DIR
archive get -C {directory in $ARC} {filename}
#
#Run the analysis.
#
{My Program}
#
#Archive output and remove $WORK_DIR
#
tar cvf ../{output filename}.tar .
archive put -C {directory in $ARC} ../{output filename}.tar
rm -rf $WORK_DIR
#
#Exit the script.
#
exit

```

This script copies an input file and a program to the user's \$WORK_DIR directory. The \$WORK_DIR directory is a directory created by LSF for users to run their batch jobs. This directory has certain protections from the workspace scrubber, as long as the job is running, plus five days after it finishes, this directory will not be removed. Then the script changes to the \$WORK_DIR directory, runs the program, copies two output files to permanent storage (one to the \$HOME directory, one to the archival storage system \$ARC), and then deletes the remaining files, the \$WORK_DIR directory and exits.

NOTE: \$WORK_DIR only exists in LSF, you will not be able to change to that directory using the variable \$WORK_DIR.

4.2.3 Submitting Jobs

Once a job script is prepared, the `bsub` command is used to submit the script to LSF. The command has the following syntax:

bsub < script

Some important LSF options used are (type `man bsub` for a complete list of options available):

- | | |
|------------------------|--|
| -q <i>queue</i> | Specifies the name of the queue to which the job will be submitted. For a list of allowable queues, please see:
http://www.asc.hpc.mil/overall/policy_procedure/policies/use_policy.php |
| -n <i>n</i> | Specifies the number of CPUs the job will use. The system schedules to the node level. As such, this number must be specified in multiples of two. |

-W hh:mm	Specifies the time limit for the job in walltime. The time should be specified in the hh:mm format (e.g., 15:00). *This is a required field, there is no default.
-o <i>outfilename</i>	Standard output (stdout) for the job is written to <i>outfilename</i> . *If you do not specify an output filename, LSF emails the output to your ASC email account.
-e <i>errfilename</i>	Standard error (stderr) for the job is written to <i>errfilename</i> . The default name is <i>jobname.ennn</i> where <i>nnn</i> is the LSF identifier. *If you do not specify an error filename, LSF emails the error information to your ASC email account.
-J jobname	Specifies the name of the job.
-P account	Specifies the account number to charge to.
-a [SMP MPI MIX]	Specifies how jobs should be spread across nodes, if at all. SMP and MIX forces all CPUs to be allocated on a single node while MPI allows CPUs to be spread across multiple nodes. This is a required field.

When a job is submitted to LSF, a unique identifier is assigned to the job by the batch system similar to below:

```
2079.falcon-1.asc.hpc.mil
```

This identifier is needed when deleting a job.

Options of `bsub` commands are specified within the script file itself. The options are specified using syntax similar to PBS, but each line that contains an option must begin with the `#BSUB` string. Options that are specified within the script file must precede the first executable shell command of the file as in the following example.

```
#!/bin/csh
#BSUB -q standard
#BSUB -n 1
#BSUB -W 168:00
#BSUB -J test
#BSUB -o test.out
#BSUB -e test.out
#BSUB -a MPI
#BSUB -P WP+WPASC00000000**
```

**This is an example number. To find your account number, check your `$ACCOUNT` variable using

echo \$ACCOUNT

More sample batch scripts can be found at the following URL:

<http://www.asc.hpc.mil/customer/userdocs/samples/samplebatch.php>

4.2.4 Monitoring Jobs

The *bjobs* command is used to report the status of the batch jobs that are currently queued or running. Type `man bjobs` for information about *bjobs* and the options that are available.

The *bjobs* command lists all jobs that are running and queued.

bjobs -u all

<u>JOBID</u>	<u>USER</u>	<u>STAT</u>	<u>QUEUE</u>	<u>FROM_HOST</u>	<u>EXEC_HOST</u>	<u>JOBNAME</u>	<u>SUBMIT_TIME</u>
3373	user1	PEND	standard	falcon1041		test	Feb 5 15:22
3971	user2	RUN	standard	falcon1042	2*ls-host.1	test	Feb 5 15:22

Here is an explanation of the fields in the *bjobs* output.

Table 2: Fields from *bjobs*

<u>Item</u>	<u>Meaning</u>
JOBID	A unique identifier that consists of the original request number and the machine from which the request was submitted. Format is <i>nnn</i> , where <i>nnn</i> is an integer.
USER	Username of person submitting the job.
STAT	Job status. “RUN” indicates the job is running; “PEND” indicates the job is queued.
QUEUE	Name of the queue where the job is waiting or executing.
FROM_HOST	Cluster domain from which the job was submitted.
EXEC_HOST	Cluster domain where the job is running.
JOBNAME	Name of the job. This is either the name of the script file submitted to LSF or the name chosen with the -J flag.
SUBMIT_TIME	The date and time the job was submitted.

4.2.5 Deleting Jobs

In LSF, queued or running jobs are removed using the `bkill` command. The syntax is

bkill request-id

where *request-id* is the LSF identifier number.

Example:

bjobs

<u>JOBID</u>	<u>USER</u>	<u>STAT</u>	<u>QUEUE</u>	<u>FROM_HOST</u>	<u>EXEC_HOST</u>	<u>JOBNAME</u>	<u>SUBMIT_TIME</u>
3373	user	PEND	standard	falcon1041	2*ls-host.1	test	Feb 5 15:22

bkill 3373

5. Customer Service

5.1 Customer Service Center

For customer assistance, call the ASC MSRC Service Center at 1-888-MSRC-ASC (1-888-677-2272), (937) 255-0194, or DSN 785-0194, or send e-mail with a description of the problem to msrchelp@asc.hpc.mil. The support analysts will help with anything related to ASC MSRC: third party software, UNIX, the different ASC MSRC computers, etc. If you have any questions about the ASC MSRC, contact the Service Center first. If your problem or question is beyond the scope of their expertise, they will refer you to the appropriate resource.

5.2 ASC MSRC Support

In-depth technical inquiries and problems are forwarded to the ASC MSRC Customer Assistance and Technology Center (CATC), which pursues such inquiries and problems through resolution as rapidly as possible. The ASC MSRC CATC will attempt to determine the nature of the problem, then identify and coordinate whatever resources are needed to resolve the problem.

The ASC MSRC also offers training classes, which provide an introduction to UNIX and the ASC MSRC. Intermediate and advanced classes on selected topics are also periodically announced on the Programming Environment and Training (PET) section of the ASC MSRC homepage. Topics for such classes may be requested through the Customer Service Center.

The ASC MSRC CATC is ready to support in an advisory capacity any engineer or scientist who is (or potentially is) an ASC MSRC user.

5.3 ASC MSRC Website

The ASC MSRC website is the best source for current ASC MSRC information. To access the ASC MSRC website simply access this URL: <http://www.asc.hpc.mil>.

Some of the topics found on the website include:

APPLICATIONS

Short and long descriptions of current ASC MSRC applications

<http://www.asc.hpc.mil/software/>

SYSTEMS

Information on ASC MSRC servers and Archival Storage

<http://www.asc.hpc.mil/hardware/>

CUSTOMER SERVICE

Available Customer Services

<http://www.asc.hpc.mil/customer/>

ONLINE DOCUMENTATION

Listings of the ASC MSRC User Guides are available for viewing. Instructions are given on obtaining postscript versions.

<http://www.asc.hpc.mil/customer/userdocs/>

VISUALIZATION LAB INFORMATION

Current status and other information about the Visualization Lab.

<http://www.asc.hpc.mil/sciviz/>

TRAINING

Current course offerings and schedule

<https://okc.erdhpc.mil/index.jsp>

FREQUENTLY ASKED QUESTIONS

Submit questions and read about various topics (such as “Customizing Your Environment”)

<http://www.asc.hpc.mil>

POLICIES AND PROCEDURES

The latest policies regarding usage of the ASC MSRC resources.

http://www.asc.hpc.mil/overall/policy_procedure/

Appendix A. Usage Hints

The following are tips and hints for the effective use of the ASC MSRC HP XC Opteron.

A.1 Runtime Considerations

A.1.1 Batch use is recommended

The ASC MSRC HP XC Opteron system allocates more resources to batch jobs than for interactive use. Users will obtain the best throughput for long running or large memory jobs by submitting jobs to the batch queues.

A.1.2 Request only the time and memory needed

When submitting a job, choose the smallest queue that accommodates the job's time and memory requirements. Jobs that request significantly more resources than are actually needed can result in longer wait times and inefficient use of the machine.

A.2 Files and Filespace

A.2.1 File Management in workspace

A file scrubber is used to automatically remove old files from workspace to prevent them from becoming filled. The policy for removing files from these filesystems is available on our website. However, you are encouraged to remove files from workspace when they are no longer needed. This will minimize the overhead needed to enforce this policy.

The workspace filesystem is not backed up. It is your responsibility to transfer files that need to be saved to a location that allows permanent storage. Two possibilities are your \$HOME directory or \$ARC. Due to space restrictions on your \$HOME directory, it is highly recommended that you use \$ARC for long-term file storage and backups.

A.2.2 Archival Storage

To provide long-term storage and archiving, the ASC MSRC provides an archival storage system that combines a large-capacity Serial ATA drive cache, local tape backup/storage and a remote disaster recovery site.

This area is referenced using the environment variable \$ARC, is NFS mounted to the interactive node of eagle and can be accessed much like any other directory in a filesystem. However, to improve system performance and increase transfer speeds, \$ARC is not NFS mounted to the batch nodes of eagle. To transfer files in the batch environment, you will need to use unkerberized rcp or the archive command.

```
/usr/bin/rcp ${msas}:/msas*/foo/bar $WRK/username
```

```
archive get archive_filename local_filename
```

```
archive put local_filename
```

For more details about the Archival Storage system, see the *Archival Storage User's Guide*, located at:

<http://www.asc.hpc.mil/customer/userdocs/>

A.2.3 Keep I/O local to the system

The workspace filesystems are local to the ASC MSRC HP XC Opteron via the lustre filesystem. Although \$HOME is NFS-mounted internally to the compute nodes, I/O access from workspace will be faster than from the HAFS.

Here is a sample script that copies two input files (one from the \$HOME directory, one from the archival storage system) and a program to the user's workspace area, changes to workspace, runs the program, copies two output files (one to the \$HOME directory, one to the archival storage system) and then deletes the remaining files.

```
cd $WRK
cp $HOME/small.input $WRK
archive get big.input $WRK
archive get prog $WRK
prog
archive put big.output
mv small.output $HOME
rm small.input big.input prog big.output
```